

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121522\
 Data File : BM038026.D
 Acq On : 15 Dec 2022 11:32
 Operator : CG/JU
 Sample : SST0.807
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8018

Quant Time: Dec 15 21:43:10 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM121522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 15 21:40:10 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.064	152	3304	0.400	ng/u1	0.00
4) Naphthalene-d8	10.878	136	9451	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.685	164	5132	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.422	188	10426	0.400	ng/u1	0.00
17) Chrysene-d12	21.583	240	7595	0.400	ng/u1	0.00
23) Perylene-d12	24.041	264	7124	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	3547	0.761	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.456	152	11565	0.864	ng/u1	0.00
18) Fluoranthene-d10	19.436	212	22370	0.863	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.448	88	3423	0.743	ng/u1	92
5) Naphthalene	10.928	128	23954	0.830	ng/u1	99
7) 2-Methylnaphthalene	12.528	142	14831	0.871	ng/u1	99
8) 1-Methylnaphthalene	12.748	142	15090	0.865	ng/u1	100
10) Acenaphthylene	14.407	152	22632	0.777	ng/u1	99
11) Acenaphthene	14.749	153	17114	0.820	ng/u1	100
12) Fluorene	15.730	166	19203	0.802	ng/u1	100
14) Pentachlorophenol	17.071	266	3538	0.798	ng/u1	99
15) Phenanthrene	17.464	178	29514	0.817	ng/u1	99
16) Anthracene	17.552	178	27622	0.836	ng/u1	99
19) Fluoranthene	19.469	202	32091	0.800	ng/u1	100
20) Pyrene	19.831	202	32621	0.801	ng/u1	100
21) Benzo(a)anthracene	21.568	228	26752	0.771	ng/u1	98
22) Chrysene	21.624	228	27259	0.757	ng/u1	100
24) Benzo(b)fluoranthene	23.284	252	28023	0.709	ng/u1	93
25) Benzo(k)fluoranthene	23.333	252	27308	0.737	ng/u1	93
26) Benzo(a)pyrene	23.930	252	24055	0.750	ng/u1#	92
27) Indeno(1,2,3-cd)pyrene	26.606	276	30046	0.770	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.622	278	23492	0.783	ng/u1	96
29) Benzo(g,h,i)perylene	27.400	276	25532	0.787	ng/u1	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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