

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111524\
 Data File : BN035101.D
 Acq On : 15 Nov 2024 12:13
 Operator : RC/JU
 Sample : P4803-07
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A0AQ2

Quant Time: Nov 15 12:42:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 00:41:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.550	152	2602	0.400	ng/ul	0.00
4) Naphthalene-d8	10.311	136	6677	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.189	164	5291	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	12329	0.400	ng/ul	0.00
17) Chrysene-d12	21.137	240	10005	0.400	ng/ul	# 0.00
23) Perylene-d12	23.299	264	9261	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1444	0.583	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.912	152	4062	0.397	ng/ul	0.00
18) Fluoranthene-d10	18.972	212	12613	0.491	ng/ul	0.00
Target Compounds				Qvalue		
5) Naphthalene	10.361	128	41026	2.455	ng/ul	96
10) Acenaphthylene	13.903	152	303060	14.020	ng/ul	98
11) Acenaphthene	14.250	153	207157	12.642	ng/ul	97
12) Fluorene	15.240	166	46746	2.242	ng/ul	99
15) Phenanthrene	16.979	178	143659	4.361	ng/ul	97
16) Anthracene	17.068	178	137714	4.448	ng/ul	97
19) Fluoranthene	19.000	202	144355	4.318	ng/ul#	96
20) Pyrene	19.362	202	147531	4.256	ng/ul#	96
21) Benzo(a)anthracene	21.119	228	67956	1.875	ng/ul	99
22) Chrysene	21.172	228	46173	1.249	ng/ul	100
24) Benzo(b)fluoranthene	22.656	252	149012	4.576	ng/ul	95
25) Benzo(k)fluoranthene	22.697	252	66234	1.943	ng/ul	95
26) Benzo(a)pyrene	23.206	252	53770	1.848	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.450	276	51078	1.435	ng/ul#	89
28) Dibenzo(a,h)anthracene	25.463	278	52042	1.877	ng/ul	95
29) Benzo(g,h,i)perylene	26.097	276	44662	1.643	ng/ul	97

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

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