

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082824\
 Data File : BN033647.D
 Acq On : 28 Aug 2024 15:57
 Operator : MA/JU
 Sample : P3711-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E24A1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 08/29/2024
 Supervised By :mohammad ahmed 08/30/2024

Quant Time: Aug 28 16:30:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 23:48:09 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.536	152	5405	0.400	ng/ul	0.00
4) Naphthalene-d8	10.299	136	14924	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.179	164	7906	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.927	188	16669m	0.400	ng/ul	0.00
17) Chrysene-d12	21.141	240	12920m	0.400	ng/ul	0.00
23) Perylene-d12	23.307	264	13450m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	26041	4.686	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.894	152	5986	0.280	ng/ul	0.00
18) Fluoranthene-d10	18.966	212	17018	0.442	ng/ul	0.00
Target Compounds						Qvalue
14) Pentachlorophenol	16.581	266	227m	0.043	ng/ul	
15) Phenanthrene	16.969	178	1079m	0.022	ng/ul	
19) Fluoranthene	18.994	202	1408	0.028	ng/ul#	75
24) Benzo(b)fluoranthene	22.664	252	1165	0.022	ng/ul#	1
25) Benzo(k)fluoranthene	22.708	252	1106	0.021	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	25.455	276	1503	0.024	ng/ul	90
28) Dibenzo(a,h)anthracene	25.472	278	1110	0.023	ng/ul#	11
29) Benzo(g,h,i)perylene	26.102	276	1268	0.026	ng/ul#	50

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

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