

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012258.D
 Acq On : 21 Oct 2022 22:24
 Operator : CG/JU
 Sample : N5064-11
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BR6

Quant Time: Oct 22 01:13:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	422202	20.000	ng/u1	0.00
20) Naphthalene-d8	10.628	136	1675866	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.469	164	952430	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.233	188	1920393	20.000	ng/u1	0.00
79) Chrysene-d12	21.321	240	1992448	20.000	ng/u1	-0.01
88) Perylene-d12	23.727	264	2133352	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	38204	3.662	ng/uL	0.00
4) Pyridine-d5	3.693	84	59398	1.976	ng/u1	0.01
7) Phenol-d5	6.993	99	660934	18.133	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	416205	18.828	ng/u1	0.00
11) 2-Chlorophenol-d4	7.351	132	540295	19.474	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	486063	17.158	ng/u1	-0.01
21) Nitrobenzene-d5	8.992	128	254390	20.507	ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	275140	20.558	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	465633	18.548	ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	624980	17.522	ng/u1	-0.01
46) Dimethylphthalate-d6	13.886	166	1351510	20.399	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1569302	20.751	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.692	143	180466	14.540	ng/u1	0.00
60) Fluorene-d10	15.469	176	1176198	20.636	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	111725	10.227	ng/u1	0.00
73) Anthracene-d10	17.333	188	1801085	21.397	ng/u1	0.00
81) Pyrene-d10	19.568	212	2168474	19.116	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.568	264	2301455	21.512	ng/u1	-0.02

Target Compounds	Qvalue

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

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