

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121123\
 Data File : BP018735.D
 Acq On : 11 Dec 2023 19:59
 Operator : MA/JU
 Sample : 04929-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AQ3

Quant Time: Dec 11 22:02:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 11 22:02:13 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	294552	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	826064	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	521277	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.222	188	1229174	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1370092	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1635821	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	37345	4.320	ng/uL	0.00
4) Pyridine-d5	3.681	84	483559	20.886	ng/ul	0.00
7) Phenol-d5	7.010	99	633712	21.439	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	389520	22.234	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	501083	21.725	ng/ul	0.00
15) 4-Methylphenol-d8	8.552	113	490167	20.490	ng/ul	0.00
21) Nitrobenzene-d5	9.004	128	231063	31.428	ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	225616	29.598	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	357027	22.906	ng/ul	0.00
31) 4-Chloroaniline-d4	10.775	131	468571	23.003	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166	1141723	24.540	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1241206	24.228	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	171851	22.503	ng/ul	0.00
60) Fluorene-d10	15.463	176	985991	25.248	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	145966	19.805	ng/ul	0.00
73) Anthracene-d10	17.316	188	1594036	24.668	ng/ul	0.00
81) Pyrene-d10	19.557	212	2062822	19.342	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	2311324	24.233	ng/ul	0.00

Target Compounds Qvalue

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

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