

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018359.D
 Acq On : 25 Nov 2023 09:42
 Operator : MA/JU
 Sample : 05264-17
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKS8

Quant Time: Nov 26 23:58:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 24 22:02:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	378194	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	1401402	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	782717	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1629021	20.000	ng/ul	0.00
79) Chrysene-d12	21.356	240	1723495	20.000	ng/ul	0.00
88) Perylene-d12	23.792	264	2050137	20.000	ng/ul	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	37770	3.741	ng/uL	0.00
4) Pyridine-d5	3.687	84	203519	7.334	ng/ul	0.00
7) Phenol-d5	7.028	99	115589	3.344	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.198	67	344556	16.671	ng/ul	0.00
11) 2-Chlorophenol-d4	7.393	132	375018	12.923	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	207595	7.443	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	235637	20.701	ng/ul	0.00
24) 2-Nitrophenol-d4	9.745	143	216499	20.595	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.286	165	401919	15.847	ng/ul	0.00
31) 4-Chloroaniline-d4	10.804	131	587544	17.177	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166	1327790	19.065	ng/ul	0.00
49) Acenaphthylene-d8	14.192	160	1521851	19.177	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	34686	3.608	ng/ul	0.00
60) Fluorene-d10	15.492	176	1139234	19.556	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	114313	16.484	ng/ul	0.00
73) Anthracene-d10	17.345	188	1620860	18.616	ng/ul	0.00
81) Pyrene-d10	19.586	212	2340083	17.060	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	2572267	21.642	ng/ul	0.01

Target Compounds	Qvalue

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

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