

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018292.D
 Acq On : 23 Nov 2023 00:52
 Operator : MA/JU
 Sample : 05226-21
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCG77

Quant Time: Nov 23 01:24:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 22 14:09:55 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	404960	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1698511	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	1062432	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2017345	20.000	ng/u1	0.00
79) Chrysene-d12	21.357	240	1199827	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1371996	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.276	96	45703	4.227	ng/uL	0.00
4) Pyridine-d5	3.687	84	301154	10.135	ng/u1	0.00
7) Phenol-d5	7.034	99	324634	8.772	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.211	67	783069	35.384	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	780325	25.112	ng/u1	0.00
15) 4-Methylphenol-d8	8.581	113	535553	17.931	ng/u1	0.00
21) Nitrobenzene-d5	9.034	128	499332	36.194	ng/u1	0.00
24) 2-Nitrophenol-d4	9.758	143	444715	34.904	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	881740	28.685	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	1303409	31.439	ng/u1	0.00
46) Dimethylphthalate-d6	13.922	166	3406688	36.036	ng/u1	-0.01
49) Acenaphthylene-d8	14.198	160	3719779	34.533	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	79024	6.055	ng/u1	-0.01
60) Fluorene-d10	15.498	176	2816249	35.616	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	255002	29.694	ng/u1	0.00
73) Anthracene-d10	17.357	188	4167250	38.648	ng/u1	0.00
81) Pyrene-d10	19.592	212	3921246	41.064	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.621	264	3238418	40.713	ng/u1	0.00

Target Compounds Qvalue

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

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