

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018288.D
 Acq On : 22 Nov 2023 22:36
 Operator : MA/JU
 Sample : 05226-13
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCG43

Quant Time: Nov 22 23:20:17 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	393014	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1661830	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.510	164	1066673	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2107057	20.000	ng/u1	0.00
79) Chrysene-d12	21.357	240	1284154	20.000	ng/u1	0.00
88) Perylene-d12	23.780	264	1358603	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	41636	3.968	ng/uL	0.00
4) Pyridine-d5	3.687	84	328924	11.407	ng/u1	0.00
7) Phenol-d5	7.034	99	218058	6.071	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	729990	33.988	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	695665	23.068	ng/u1	0.00
15) 4-Methylphenol-d8	8.581	113	461107	15.908	ng/u1	0.00
21) Nitrobenzene-d5	9.034	128	467722	34.651	ng/u1	0.00
24) 2-Nitrophenol-d4	9.757	143	416638	33.422	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	828097	27.534	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	1194788	29.455	ng/u1	0.00
46) Dimethylphthalate-d6	13.928	166	3221983	33.947	ng/u1	0.00
49) Acenaphthylene-d8	14.204	160	3593852	33.231	ng/u1	0.00
54) 4-Nitrophenol-d4	14.687	143	52116	3.977	ng/u1	-0.01
60) Fluorene-d10	15.504	176	2691334	33.901	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	250649	27.944	ng/u1	0.00
73) Anthracene-d10	17.357	188	4048708	35.950	ng/u1	0.00
81) Pyrene-d10	19.592	212	3929686	38.450	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.621	264	2912264	36.974	ng/u1	0.00

Target Compounds Qvalue

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

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