

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112423\
 Data File : BP018326.D
 Acq On : 24 Nov 2023 14:29
 Operator : MA/JU
 Sample : 05268-21
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCKB6

Quant Time: Nov 24 21:47:29 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	7.863	152	428700	20.000	ng/u1	-0.02
20) Naphthalene-d8	10.663	136	1666895	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.504	164	920711	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.251	188	1640588	20.000	ng/u1	-0.01
79) Chrysene-d12	21.363	240	1412524	20.000	ng/u1	0.00
88) Perylene-d12	23.792	264	1821954	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.264	96	42390	3.704	ng/uL	-0.01
4) Pyridine-d5	3.681	84	144030	4.579	ng/u1	0.00
7) Phenol-d5	7.016	99	173039	4.417	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.193	67	577194	24.637	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.387	132	586047	17.816	ng/u1	-0.02
15) 4-Methylphenol-d8	8.563	113	342579	10.835	ng/u1	-0.02
21) Nitrobenzene-d5	9.022	128	375389	27.726	ng/u1	-0.02
24) 2-Nitrophenol-d4	9.746	143	329811	26.377	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.281	165	654803	21.706	ng/u1	-0.02
31) 4-Chloroaniline-d4	10.799	131	798584	19.628	ng/u1	-0.02
46) Dimethylphthalate-d6	13.916	166	2218607	27.081	ng/u1	-0.02
49) Acenaphthylene-d8	14.193	160	2486825	26.640	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.687	143	36885	3.261	ng/u1	-0.01
60) Fluorene-d10	15.498	176	1862233	27.176	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	168722	24.159	ng/u1	-0.01
73) Anthracene-d10	17.351	188	2787031	31.783	ng/u1	-0.01
81) Pyrene-d10	19.592	212	2946025	26.206	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.633	264	3342794	31.647	ng/u1	0.00

Target Compounds Qvalue

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

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