

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112223\
 Data File : BP018297.D
 Acq On : 23 Nov 2023 03:41
 Operator : MA/JU
 Sample : PB156976BL
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK976

Quant Time: Nov 23 04:12:37 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	404970	20.000	ng/u1	0.00
20) Naphthalene-d8	10.675	136	1732597	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.504	164	1155945	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.257	188	2385912	20.000	ng/u1	0.00
79) Chrysene-d12	21.357	240	1462868	20.000	ng/u1	0.00
88) Perylene-d12	23.774	264	1433476	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	62580	5.788	ng/uL	0.00
4) Pyridine-d5	3.687	84	880454	29.631	ng/u1	0.00
7) Phenol-d5	7.034	99	1066039	28.804	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	658039	29.733	ng/u1	0.00
11) 2-Chlorophenol-d4	7.405	132	878328	28.265	ng/u1	0.00
15) 4-Methylphenol-d8	8.581	113	881949	29.528	ng/u1	0.00
21) Nitrobenzene-d5	9.034	128	433521	30.806	ng/u1	0.00
24) 2-Nitrophenol-d4	9.757	143	419322	32.264	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	872014	27.810	ng/u1	0.00
31) 4-Chloroaniline-d4	10.810	131	1341374	31.718	ng/u1	0.00
46) Dimethylphthalate-d6	13.922	166	3131254	30.443	ng/u1	-0.01
49) Acenaphthylene-d8	14.198	160	3437094	29.327	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	376585	26.521	ng/u1	0.00
60) Fluorene-d10	15.498	176	2559249	29.748	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.610	200	236431	23.278	ng/u1	0.00
73) Anthracene-d10	17.357	188	3800003	29.798	ng/u1	0.00
81) Pyrene-d10	19.592	212	3745602	32.171	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.621	264	2539639	30.559	ng/u1	0.00

Target Compounds Qvalue

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

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