

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051623\
 Data File : BP015101.D
 Acq On : 16 May 2023 09:55
 Operator : CG/JU
 Sample : PB152799BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK799

Quant Time: May 17 00:43:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 16 01:39:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	302883	20.000	ng/u1	0.00
20) Naphthalene-d8	10.963	136	1414911	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.775	164	926458	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.534	188	2049460	20.000	ng/u1	0.00
79) Chrysene-d12	21.610	240	1744389	20.000	ng/u1	0.00
88) Perylene-d12	24.186	264	1780135	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	51132	6.559	ng/uL	0.00
4) Pyridine-d5	3.870	84	733380	33.190	ng/u1	0.00
7) Phenol-d5	7.281	99	890854	31.678	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.452	67	573200	35.292	ng/u1	0.00
11) 2-Chlorophenol-d4	7.658	132	711548	34.413	ng/u1	0.00
15) 4-Methylphenol-d8	8.846	113	763714	34.081	ng/u1	0.00
21) Nitrobenzene-d5	9.310	128	368403	33.135	ng/u1	0.00
24) 2-Nitrophenol-d4	10.040	143	410032	33.144	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.581	165	713257	31.951	ng/u1	0.00
31) 4-Chloroaniline-d4	11.099	131	1132009	34.273	ng/u1	0.00
46) Dimethylphthalate-d6	14.181	166	2468867	35.265	ng/u1	0.00
49) Acenaphthylene-d8	14.475	160	2787518	34.814	ng/u1	0.00
54) 4-Nitrophenol-d4	14.963	143	413823	30.441	ng/u1	0.00
60) Fluorene-d10	15.769	176	2071071	35.059	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.887	200	348972	27.019	ng/u1	0.00
73) Anthracene-d10	17.634	188	3283040	34.983	ng/u1	0.00
81) Pyrene-d10	19.851	212	3748572	37.052	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.016	264	3299345	37.337	ng/u1	0.00

Target Compounds Qvalue

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

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