

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092923\
 Data File : BP017458.D
 Acq On : 29 Sep 2023 22:25
 Operator : MA/JU
 Sample : 04497-16
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJ22

Quant Time: Sep 30 00:39:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 29 22:31:03 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.952	152	232197	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.793	136	911156	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.645	164	535993	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.433	188	1183377	20.000	ng/u1	-0.01
79) Chrysene-d12	21.563	240	1196900	20.000	ng/u1	-0.01
88) Perylene-d12	24.133	264	1383488	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	16602	2.937	ng/uL	0.00
4) Pyridine-d5	3.734	84	160016	10.125	ng/u1	0.00
7) Phenol-d5	7.105	99	126720	6.645	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.293	67	316850	27.531	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.469	132	345649	22.853	ng/u1	-0.01
15) 4-Methylphenol-d8	8.669	113	221640	14.949	ng/u1	-0.02
21) Nitrobenzene-d5	9.157	128	202519	30.745	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.875	143	213246	29.815	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.404	165	366505	27.123	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.957	131	462155	23.610	ng/u1	-0.01
46) Dimethylphthalate-d6	14.057	166	1100832	28.670	ng/u1	-0.01
49) Acenaphthylene-d8	14.340	160	1281511	29.019	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.881	143	32620	4.876	ng/u1	-0.01
60) Fluorene-d10	15.645	176	967857	29.279	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.787	200	138734	20.800	ng/u1	-0.02
73) Anthracene-d10	17.533	188	1584568	30.192	ng/u1	-0.01
81) Pyrene-d10	19.780	212	1952886	30.172	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.968	264	2071478	31.098	ng/u1	-0.02

Target Compounds Qvalue

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

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