

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020325\
 Data File : BP023884.D
 Acq On : 03 Feb 2025 17:55
 Operator : RC/JU
 Sample : PB166466BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK466

Quant Time: Feb 04 07:33:45 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Feb 03 11:19:36 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	362748	20.000	ng/u1	0.00
20) Naphthalene-d8	10.557	136	1560330	20.000	ng/u1	0.01
38) Acenaphthene-d10	14.404	164	1002862	20.000	ng/u1	0.02
64) Phenanthrene-d10	17.198	188	2054338	20.000	ng/u1	0.01
79) Chrysene-d12	21.633	240	2008701	20.000	ng/u1	0.00
88) Perylene-d12	25.004	264	2187980	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	59800	6.849	ng/uL	0.00
4) Pyridine-d5	3.711	84	885198	35.137	ng/u1	0.00
7) Phenol-d5	6.964	99	1145383	35.679	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.117	67	586031	34.488	ng/u1	0.01
11) 2-Chlorophenol-d4	7.328	132	908067	36.022	ng/u1	0.01
15) 4-Methylphenol-d8	8.481	113	901843	34.390	ng/u1	0.00
21) Nitrobenzene-d5	8.922	128	410744	32.903	ng/u1	0.00
24) 2-Nitrophenol-d4	9.640	143	420959	31.251	ng/u1	0.01
28) 2,4-Dichlorophenol-d3	10.181	165	833422	33.727	ng/u1	0.01
31) 4-Chloroaniline-d4	10.687	131	1207085	32.663	ng/u1	0.01
46) Dimethylphthalate-d6	13.804	166	2505237	33.491	ng/u1	0.01
49) Acenaphthylene-d8	14.093	160	2913220	34.530	ng/u1	0.01
54) 4-Nitrophenol-d4	14.610	143	429820	28.933	ng/u1	0.00
60) Fluorene-d10	15.410	176	2195016	34.845	ng/u1	0.02
65) 4,6-Dinitro-2-methylph...	15.528	200	267711	21.123	ng/u1	0.02
73) Anthracene-d10	17.298	188	3536032	35.056	ng/u1	0.01
81) Pyrene-d10	19.663	212	3974774	36.930	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.786	264	4050083	35.478	ng/u1	0.01

Target Compounds Qvalue

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

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